

<u>NEWS</u>	<u>1</u>	Web Page URLs for STN Seminar Schedule - N. America
<u>NEWS</u>	<u>2</u>	"Ask CAS" for self-help around the clock
<u>NEWS</u>	<u>3</u>	May 12 EXTEND option available in structure searching
<u>NEWS</u>	<u>4</u>	May 12 Polymer links for the POLYLINK command completed in REGISTRY
<u>NEWS</u>	<u>5</u>	May 27 New UPM (Update Code Maximum) field for more efficient patent SDIs in CAplus
<u>NEWS</u>	<u>6</u>	May 27 CAplus super roles and document types searchable in REGISTRY
<u>NEWS</u>	<u>7</u>	Jun 22 STN Patent Forums to be held July 19-22, 2004
<u>NEWS</u>	<u>8</u>	Jun 28 Additional enzyme-catalyzed reactions added to CASREACT
<u>NEWS</u>	<u>9</u>	Jun 28 ANTE, AQUALINE, BIOENG, CIVILENG, ENVIROENG, MECHENG, and WATER from CSA now available on STN(R)
<u>NEWS EXPRESS</u>		MARCH 31 CURRENT WINDOWS VERSION IS V7.00A, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 26 APRIL 2004
<u>NEWS HOURS</u>		STN Operating Hours Plus Help Desk Availability
<u>NEWS INTER</u>		General Internet Information
<u>NEWS LOGIN</u>		Welcome Banner and News Items
<u>NEWS PHONE</u>		Direct Dial and Telecommunication Network Access to STN
<u>NEWS WWW</u>		CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

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FILE 'HOME' ENTERED AT 03:56:14 ON 08 JUL 2004

FILE 'REGISTRY' ENTERED AT 03:56:21 ON 08 JUL 2004  
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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 6 JUL 2004 HIGHEST RN 705249-96-3  
DICTIONARY FILE UPDATES: 6 JUL 2004 HIGHEST RN 705249-96-3

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2004

Please note that search-term pricing does apply when conducting SmartSELECT searches.

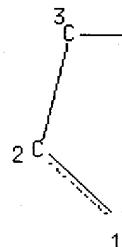
Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more

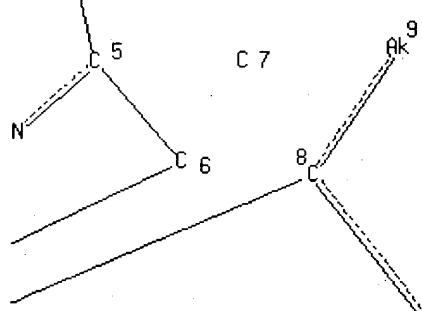
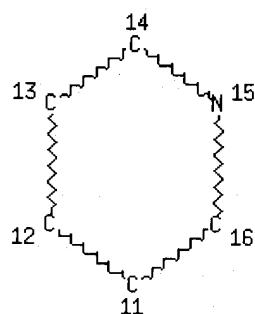
information enter **HELP\_PROP** at an arrow prompt in the file or refer to the file summary sheet on the web at:  
<http://www.cas.org/ONLINE/DBSS/registryss.html>

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L1       STRUCTURE UPLOADED

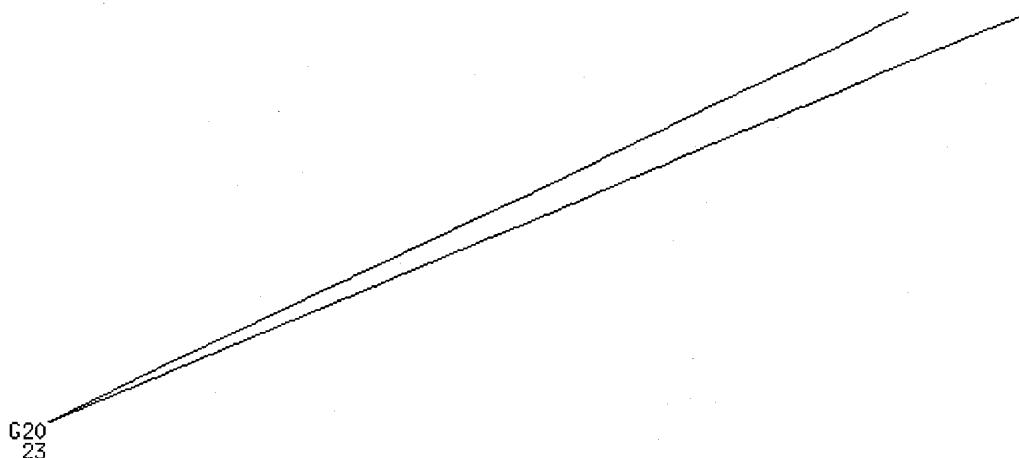
=> d 11  
L1 HAS NO ANSWERS  
L1           STR



Page 1-A

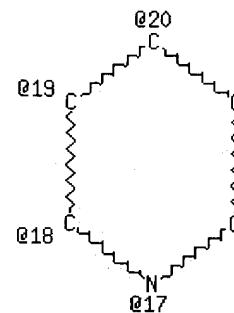


Page 1-B



Page 2-A

Hk  
@10



Page 2-B

@21

@22

Page 2-C

REP G20=(0-2) 7-6 7-8  
VPA 10-17/18/19/20/21/22 S

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MLEVEL IS CLASS AT 6 7 8 9 10
DEFAULT ECLEVEL IS LIMITED

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## GRAPH ATTRIBUTES:

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RSPEC I
NUMBER OF NODES IS 23

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## STEREO ATTRIBUTES: NONE

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SAMPLE SCREEN SEARCH COMPLETED - 6135 TO ITERATE

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16.3% PROCESSED 1000 ITERATIONS 0 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

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FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 118005 TO 127395
PROJECTED ANSWERS: 0 TO 0

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L2 0 SEA SSS SAM L1
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DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N or END:y
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SEARCH TIME: 00.00.08

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L3 4 SEA SSS FUL L1
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=> file hcaplus
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                           ENTRY SESSION
FULL ESTIMATED COST           158.36 158.57

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FILE 'HCAPLUS' ENTERED AT 04:01:22 ON 08 JUL 2004
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FILE COVERS 1907 - 8 Jul 2004 VOL 141 ISS 2  
FILE LAST UPDATED: 7 Jul 2004 (20040707/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

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=> s 13
L4          2 L3

=> s 14 and mullican, m?/au
      57 MULLICAN, M?/AU
L5          0 L4 AND MULLICAN, M?/AU

=> s 14 and lauffer, d?/au
      46 LAUFFER, D?/AU
L6          0 L4 AND LAUFFER, D?/AU

=> s 14 and tung, r?/au
      302 TUNG, R?/AU
L7          0 L4 AND TUNG, R?/AU

=> d 14, ibib abs fhitstr, 1-2
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L4 ANSWER 1 OF 2 HCAPLUS COPYRIGHT 2004 ACS on STN

Full  Citing  
 Text  References

ACCESSION NUMBER: 2001:185778 HCAPLUS  
 DOCUMENT NUMBER: 134:237837  
 TITLE: Preparation of dolastatin peptides  
 INVENTOR(S): Petit, George R.; Srirangam, Jayaram K.; Williams, Michael D.; Durkin, Kieran P. M.; Barlozzari, Teresa; Kling, Andreas; Janssen, Bernd; Haupt, Andreas  
 PATENT ASSIGNEE(S): Basf Aktiengesellschaft, Germany; Arizona Board of Regents  
 SOURCE: PCT Int. Appl., 55 pp.  
 CODEN: PIXXD2  
 DOCUMENT TYPE: Patent  
 LANGUAGE: English  
 FAMILY ACC. NUM. COUNT: 1  
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2001018032	A2	20010315	WO 2000-US24658	20000908
WO 2001018032	A3	20020711		
W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM			
RW:	GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY,			

DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ,  
CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

US 6323315

B1 20011127

US 2000-539935

20000331

PRIORITY APPLN. INFO.:

US 1999-394962

A1 19990910

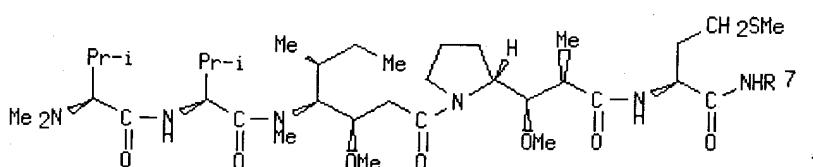
US 2000-539935

A1 20000331

OTHER SOURCE(S):

MARPAT 134:237837

GI



AB Peptides R1R2NCHR3CONHCHR4CONMeCH(R5)CH(OMe)CH2CONRCH(R6)CH(R7)  
[R2 = (CH<sub>2</sub>)<sub>3</sub>; R1-R5 are each independently H or C<sub>1</sub>-C<sub>6</sub> alkyl; A is a  
methionyl, phenylalanyl or phenylglycyl residue; n is 0 or 1; R6 is H and  
R7 is a carbocyclic, arom., alkyl, pyridylalkyl, or heterocyclic group or  
R6 is benzyl or carbalkoxy and R7 is 2-thiazolyl] or their  
pharmaceutically acceptable salts were prepd. for use as cell growth  
inhibitors. Thus, I [R7 = bicyclo[3.3.0]octan-1-yl] was prepd. by soln.  
phase methods and evaluated for in vitro cytotoxicity against a panel of  
cultured cancer cell lines, including OVCAR-3 (ovarian cancer), ED<sub>50</sub> = 3.1  
x 10<sup>-4</sup> µg/mL.

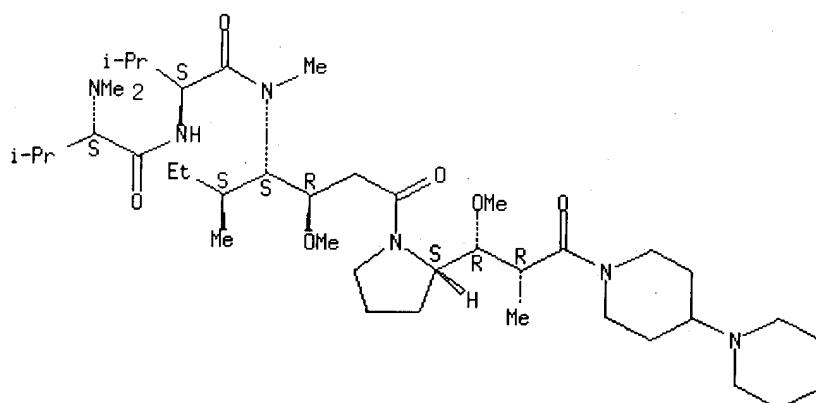
IT 329792-17-8P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological  
study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use);  
BIOL (Biological study); PREP (Preparation); USES (Uses)  
(prepn. of dolastatin peptides)

RN 329792-17-8 HCPLUS

CN L-Valinamide, N,N-dimethyl-L-valyl-N-[(1S,2R)-4-[(2S)-2-[(1R,2R)-3-[1,4'-  
bipiperidin]-1'-yl-1-methoxy-2-methyl-3-oxopropyl]-1-pyrrolidinyl]-2-  
methoxy-1-[(1S)-1-methylpropyl]-4-oxobutyl]-N-methyl- (9CI) (CA INDEX  
NAME)

Absolute stereochemistry. Rotation (-).

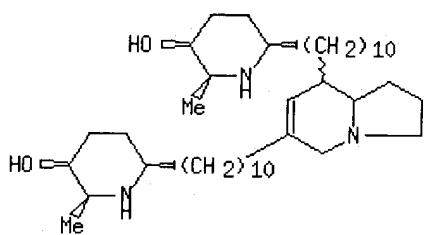


L4 ANSWER 2 OF 2 HCPLUS COPYRIGHT 2004 ACS on STN

Full Text	Citing References
-----------	-------------------

ACCESSION NUMBER: 1981:44042 HCPLUS  
DOCUMENT NUMBER: 94:44042

TITLE: Organic natural products. Part 176. Structure of the  
 alkaloid juliprosopine from Prosopis juliflora A. DC  
 AUTHOR(S): Ott-Longoni, Rita; Viswanathan, Narayananaiyer; Hesse,  
 Manfred  
 CORPORATE SOURCE: Org.-Chem. Inst., Univ. Zurich, Zurich, CH-8057,  
 Switz.  
 SOURCE: Helvetica Chimica Acta (1980), 63(7), 2119-29  
 DOCUMENT TYPE: Journal  
 LANGUAGE: German  
 GI



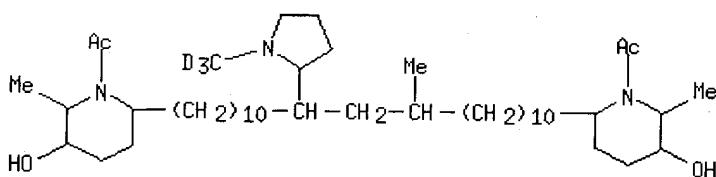
**AB** A new alkaloid, juliprosopine (I) was isolated from *P. juliflora* and its structure elucidated. The piperidine moiety could be elucidated by spectroscopic data (mass spectroscopy,  $^1\text{H-NMR}$ ,  $^{13}\text{C-NMR}$ ) of I and various derivs. The central hexahydroindolizine part and its substitution pattern were detd. mainly by the Hofmann-degrdn. product. I is built up in nature by 2 monomeric *Prosopis* alkaloids and 1 dihydropyrrol unit.

**IT 76202-04-5P**

RL: SPN (Synthetic preparation); PREP (Preparation)  
(prepn. of)

RN 76202-04-5 HCPLUS

CN 3-Piperidinol, 6,6'-(11-methyl-13-[1-(methyl-d3)-2-pyrrolidinyl]-1,23-tricosanediyl)bis[1-acetyl-2-methyl- (9CI) (CA INDEX NAME)



=> file caold

COST IN U.S. DOLLARS

SINCE FILE

ENTRY

TOTAL

SESSION

FULL ESTIMATED COST

17.42

175.99

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

ENTRY

TOTAL

CA SUBSCRIBER PRICE

-1.47

-1.47

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FILE COVERS 1907-1966  
 FILE LAST UPDATED: 01 May 1997 (19970501/UP)

This file contains CAS Registry Numbers for easy and accurate substance identification. Title keywords, authors, patent assignees, and patent information, e.g., patent numbers, are now searchable from 1907-1966. TIFF images of CA abstracts printed between 1907-1966 are available in the PAGE display formats.

This file supports REGISTRY for direct browsing and searching of all substance data from the REGISTRY file. Enter HELP FIRST for more information.

=> d his

(FILE 'HOME' ENTERED AT 03:56:14 ON 08 JUL 2004)

FILE 'REGISTRY' ENTERED AT 03:56:21 ON 08 JUL 2004

L1 STRUCTURE uploaded  
 L2 0 S L1  
 L3 4 S L1 FULL

FILE 'HCAPLUS' ENTERED AT 04:01:22 ON 08 JUL 2004

L4 2 S L3  
 L5 0 S L4 AND MULLICAN, M?/AU  
 L6 0 S L4 AND LAUFFER, D?/AU  
 L7 0 S L4 AND TUNG, R?/AU

FILE 'CAOLD' ENTERED AT 04:03:14 ON 08 JUL 2004

=> s 13  
 L8 0 L3

=> file bielstein  
 'BIELSTEIN' IS NOT A VALID FILE NAME  
 SESSION CONTINUES IN FILE 'CAOLD'  
 Enter "HELP FILE NAMES" at an arrow prompt (=>) for a list of files that are available. If you have requested multiple files, you can specify a corrected file name or you can enter "IGNORE" to continue accessing the remaining file names entered.

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.42	176.41
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-1.47

FILE 'BEILSTEIN' ENTERED AT 04:03:27 ON 08 JUL 2004  
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FILE RELOADED ON OCTOBER 20, 2002  
 FILE LAST UPDATED ON JUNE 15, 2004

FILE COVERS 1771 TO 2003.  
 FILE CONTAINS 8,997,153 SUBSTANCES

>>> PLEASE NOTE: Reaction data and substance data are stored in separate documents and can not be searched together in one query.

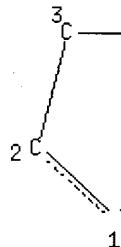
Reaction data for BEILSTEIN compounds may be displayed immediately with the display codes PRE (preparations) and REA (reactions). A substance answer set retrieved after the search for a chemical name, a molecular formula or a structure search for example can be restricted to compounds with available reaction information by concatenation with PRE/FA, REA/FA or more general with RX/FA. The BEILSTEIN Registry Number (BRN) is the link between a BEILSTEIN compound and belonging reactions. For more detailed reaction searches BRNs can be selected from substance answer sets and searched in the next step as reaction partner BRNs - Reactant (RX.RBRN) or Product BRN (RX.PBRN). After a search for reaction details substance documents associated with reactants or products may be retrieved by searching RX.PBRNs or RX.RBRNs as BRNs. <<<

>>> FOR SEARCHING PREPARATIONS SEE HELP PRE <<<

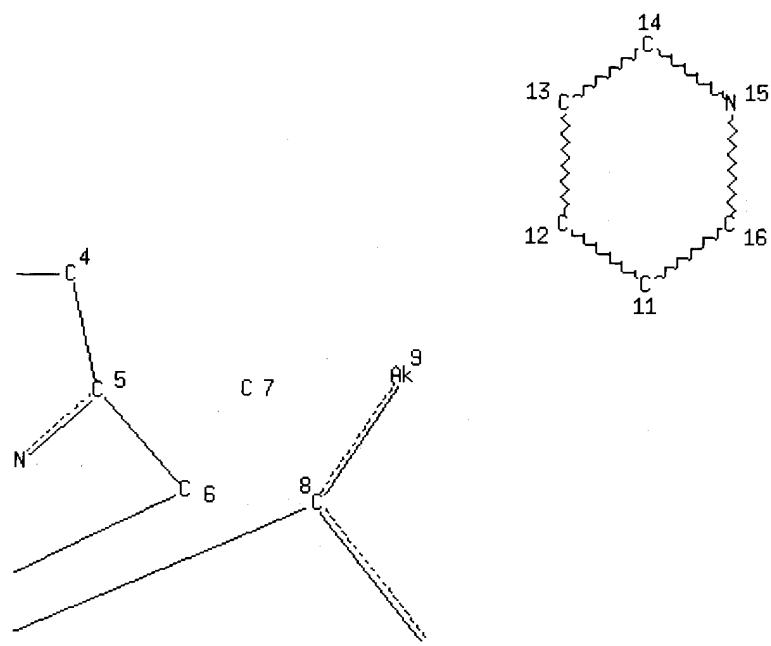
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\* PLEASE NOTE THAT THERE ARE NO FORMATS FREE OF COST. \*  
\* SET NOTICE FEATURE: THE COST ESTIMATES CALCULATED FOR SET NOTICE \*  
\* ARE BASED ON THE HIGHEST PRICE CATEGORY. THEREFORE, THESE \*  
\* ESTIMATES MAY NOT REFLECT THE ACTUAL COSTS. \*  
\* FOR PRICE INFORMATION SEE HELP COST \*  
\*\*\*\*\*

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L9        STRUCTURE UPLOADED

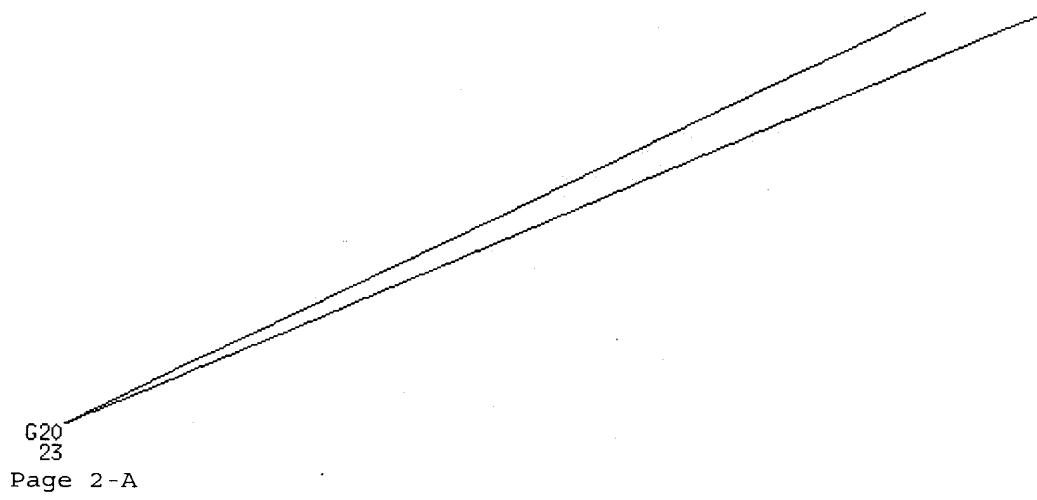
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L9 HAS NO ANSWERS  
L9        STR



Page 1-A

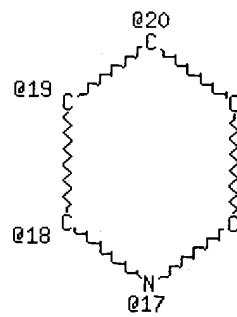


Page 1-B



Page 2-A

Ak  
@10



Page 2-B

Q21

Q22

Page 2-C

REP G20=(0-2) 7-6 7-8  
VPA 10-17/18/19/20/21/22 S

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NSPEC IS R AT 3  
NSPEC IS R AT 4  
NSPEC IS R AT 5  
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MLEVEL IS CLASS AT 6 7 8 9 10  
DEFAULT ECLEVEL IS LIMITED

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RSPEC I

NUMBER OF NODES IS 23

STEREO ATTRIBUTES: NONE

=> S 19

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SAMPLE SCREEN SEARCH COMPLETED - 285 TO ITERATE

100.0% PROCESSED 285 ITERATIONS  
SEARCH TIME: 00.00.04

0 ANSWERS

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*  
PROJECTED ITERATIONS: 4688 TO 6712  
PROJECTED ANSWERS: 0 TO 0

L10 0 SEA SSS SAM L9

=>